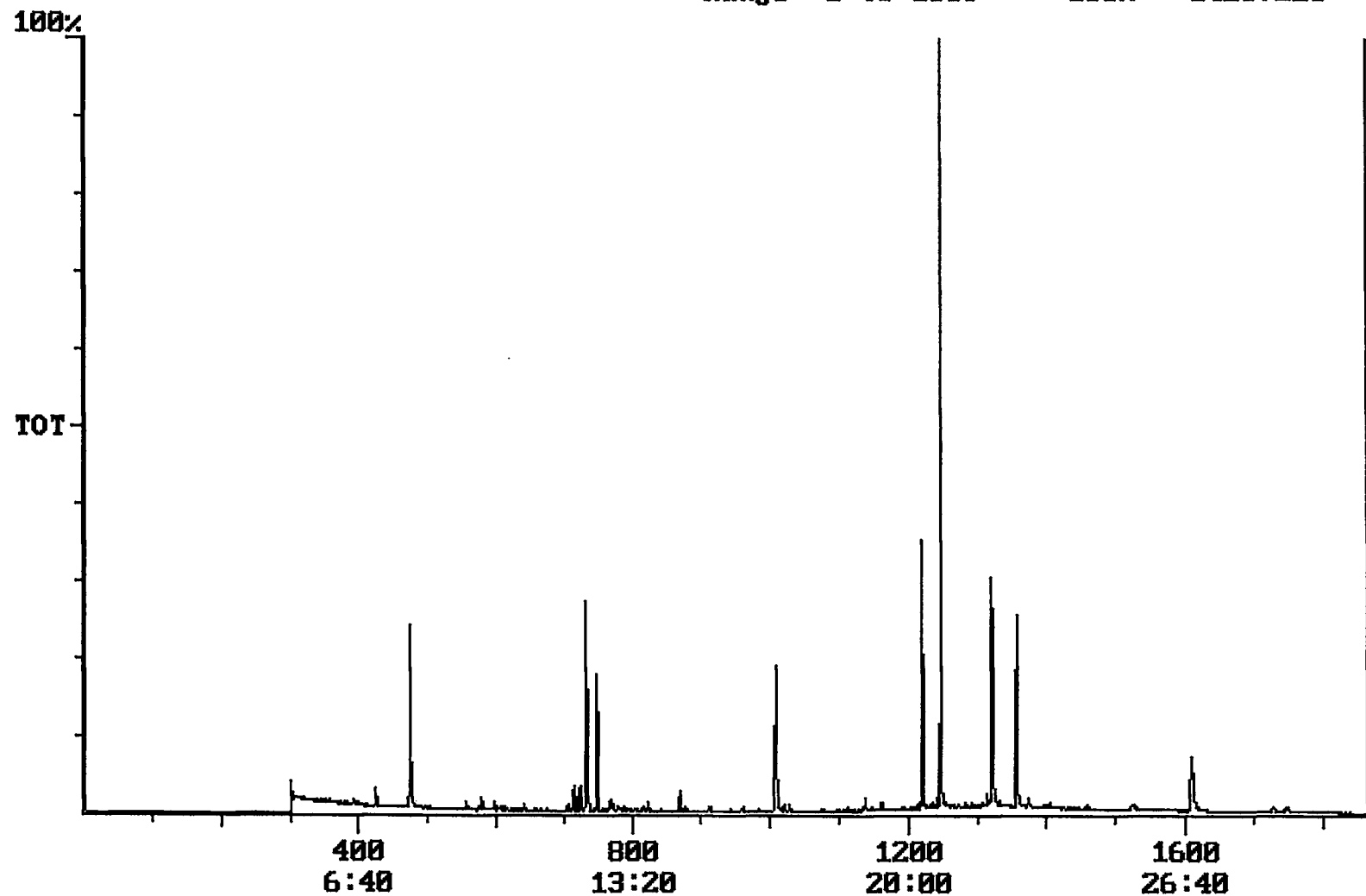


User: Bohlk, Ted
Westchester Dept. of Labs & Research
Date: 01/02/2002 Time: 09:54:10

Sample: AD26831
Analysis: \$525 Organic Chemicals - 525.2
Units: ug
Started: 11/9/01 09:52 Ended: 11/29/01 09:52 Analyst: TB
Result source: manual entry
Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	Hexachlorocyclopentadiene		< MDL		0.10
2	Hexachlorobenzene		< MDL		0.10
3	Simazine		< MDL		0.40
4	Atrazine		< MDL		0.20
5	Alachlor		< MDL		0.20
6	bis(2-Ethylhexyl)adipate		< MDL		0.60
7	bis(2-Ethylhexyl)phthalate		< MDL		0.60
8	Benzo(a)pyrene		< MDL		0.20
9	EPTC		< MDL		1.0
10	2,6-Dinitrotoluene		< MDL		2.0
11	2,4-Dinitrotoluene		< MDL		2.0
12	Molinate		< MDL		0.90
13	Terbacil		< MDL		2.0
14	Acetochlor		< MDL		2.0
15	Metribuzin		< MDL		0.15
16	Metolachlor		< MDL		0.75
17	Butachlor		< MDL		0.40
18	4,4-DDE		< MDL		0.80
19	Surr_1,3-Dimethyl-2-nitrobenze		5.2		1.0
20	Surr_Triphenyl phosphate		5.3		1.0
21	Surr_Perylene-d12		4.9		1.0

Chromatogram Plot File: E:\26831E Date: Nov-30-1991 00:06:21
Comment: BOHLK; METHOD 525; INST 204; 11/29/01; SAMPLE 26831
Scan No: 1 Retention Time: 0:01 RIC: 0 Mass Range: 0 - 0
Plotted: 1 to 1859 Range: 1 to 1859 100% = 14207220



Integration Report Quanfile: 26831E Quan Entries: 18
 Comment: BOHLK; METHOD 525; INST 204; 11/29/01; SAMPLE 26831
 Sorted via: Entry Number ↑

No	Name of Compound	R Time	Scan#	Mode	Peak Area	Pk Height
1	acenaphthene d-10 (IS)	12:13	733	Auto	1,509,855	892,563
2	phenanthrene d-10 (IS)	16:49	1009	Auto	2,413,606	876,750
3	chrysene d-12 (IS)	22:36	1356	Auto	2,069,388	1,018,269
4	p-terphenyl d-14 (IS)	20:47	1247	Auto	6,417,932	4,747,261
5	acenaphthene d-10 (SURR)	12:13	733	Auto	1,509,855	892,563
6	phenanthrene d-10 (SURR)	16:49	1009	Auto	2,413,606	876,750
7	chrysene d-12 (SURR)	22:36	1356	Auto	2,069,388	1,018,269
8	1,3-dimethyl-2-nitrobenzene	7:54	474	Auto	938,264	517,698
9	triphenyl phosphate (S)	22:01	1321	Auto	1,217,975	534,162
10	perylene d-12 (SURR)	26:49	1609	Auto	1,182,018	238,161
11	hexachlorocyclopentadiene	10:04	604	Auto	2,547	1,503
12	hexachlorobenzene	15:15	915	Auto	14,679	5,057
13	bis(2-ethylhexyl)adipate	21:54	1314	Auto	34,741	6,584
14	bis(2-ethylhexyl)phthalate	22:53	1373	Auto	144,165	51,328
15	2,6-dinitrotoluene	11:53	713	Auto	74,266	44,877
16	2,4-dinitrotoluene	13:01	781	Auto	11,168	4,827
17	butachlor	20:20	1220	Auto	1,953	1,953
18	4,4'-dichlorodiphenyl ether	20:42	1242	Auto	7,889	5,163

reversed IS / surr

Quantitation Report Quanfile: 26831E Quan Entries: 18
 Comment: BOHLK; METHOD 525; INST 204; 11/29/01; SAMPLE 26831
 Sorted via: Entry Number ↑ (S) = Standard

Cal	Name of Compound	S	Scan#	R Time	Me	Calc Amt(A)	Units
1	acenaphthene d-10(IS)	I	733	12:13	BV	5.000	UG/L
2	phenanthrene d-10 (IS)	I	1009	16:49	BV	5.000	UG/L
3	chrysene d-12(IS)	I	1356	22:36	BV	5.000	UG/L
4	p-terphenyl d-14(IS)	I	1247	20:47	BB	5.000	UG/L
5	acenaphthene d-10(SURR)	A	733	12:13	BV	1.552	UG/L
6	phenanthrene d-10(SURR)	A	1009	16:49	BV	1.727	UG/L
7	chrysene d-12 (SURR)	A	1356	22:36	BV	1.926	UG/L
8	1,3-dimethyl-2-nitrobenzene	A	474	7:54	BV	5.179	UG/L
9	triphenyl phosphate (S)	A	1321	22:01	BV	5.309	UG/L
10	perylene d-12 (SURR)	A	1609	26:49	BV	4.948	UG/L
11	hexachlorocyclopentadiene	A	604	10:04	BB	0.039	UG/L
12	hexachlorobenzene	A	915	15:15	BB	0.081	UG/L
16	bis(2-ethylhexyl)adipate	A	1314	21:54	MM	0.074	UG/L
17	bis(2-ethylhexyl)phthalate	A	1373	22:53	VB	0.186	UG/L
20	2,6-dinitrotoluene	A	713	11:53	BB	0.840	UG/L
21	2,4-dinitrotoluene	A	781	13:01	BB	0.112	UG/L
27	butachlor	A	1220	20:20	BB	0.012	UG/L
28	4,4'-dieldrin	A	1242	20:42	BB	0.040	UG/L

Comments for Sample AD26831 - Analysis \$525

Westchester Dept. of Labs & Research

User: Vinci, David L

Date: 01-03-2002 Time: 12:31:43

The recoveries for 3 of the 18 compounds in the LCS Standard were below their acceptance range. The breakdown for Endrin in the Breakdown Standard was 62%.